

n is 2 or 3;

G<sup>1</sup> is -O-, -S-, or -N(R<sup>5</sup>)-, wherein R<sup>5</sup> is -H or C<sub>1</sub>-C<sub>4</sub> alkyl; and

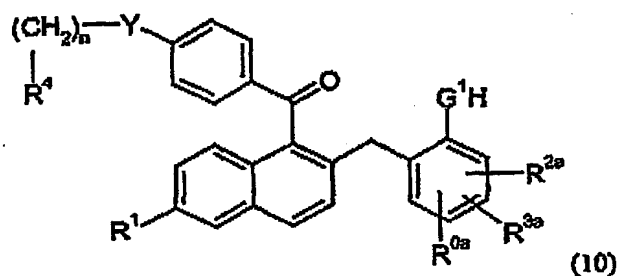
Y is -O-, -S-, -NH-, -NMe-, or -CH<sub>2</sub>-;

5 or a pharmaceutically acceptable salt thereof.

32. A compound according to Claim 31 wherein said compound is [6-hydroxy-2-(2-hydroxy-benzyl)-benzo[b]thiophen-3-yl]-[4-(2-piperidin-1-yl-ethoxy)-phenyl]-methanone.

10

33. A compound of the formula



15 wherein

R<sup>1</sup> is -H, -OH, -O(C<sub>1</sub>-C<sub>4</sub> alkyl), -OCOC<sub>6</sub>H<sub>5</sub>, -OCO(C<sub>1</sub>-C<sub>6</sub> alkyl), or -OSO<sub>2</sub>(C<sub>2</sub>-C<sub>6</sub> alkyl);

R<sup>0a</sup>, R<sup>2a</sup> and R<sup>3a</sup> are each independently -H, -OPg, or halo, wherein Pg is a hydroxy protecting group;

20 R<sup>4</sup> is 1-piperidinyl, 1-pyrrolidinyl, methyl-1-pyrrolidinyl, dimethyl-1-pyrrolidinyl, 4-morpholino, dimethylamino, diethylamino, diisopropylamino, or 1-hexamethyleneimino;

n is 2 or 3;

G<sup>1</sup> is -O-, -S-, or -N(R<sup>5</sup>)-, wherein R<sup>5</sup> is -H or C<sub>1</sub>-C<sub>4</sub> alkyl; and

25 Y is -O-, -S-, -NH-, -NMe-, or -CH<sub>2</sub>-;

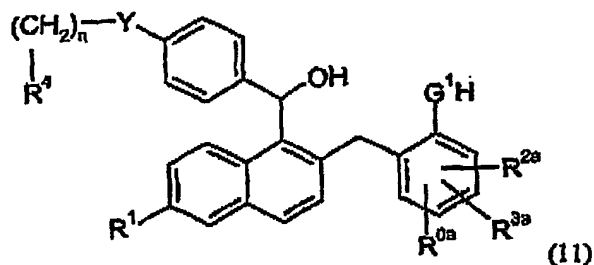
or a pharmaceutically acceptable salt thereof.

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ART 34 AMDT

-47-

34. A compound according to Claim 33 wherein said compound is [6-hydroxy-2-(2-hydroxy-benzyl)-naphthalen-1-yl]-[4-(2-piperidin-1-yl-ethoxy)-phenyl]-methanone.

5 35. A compound of the formula



wherein

10 R<sup>1</sup> is -H, -OH, -O(C<sub>1</sub>-C<sub>4</sub> alkyl), -OCOC<sub>6</sub>H<sub>5</sub>, -OCO(C<sub>1</sub>-C<sub>6</sub> alkyl), or -OSO<sub>2</sub>(C<sub>2</sub>-C<sub>6</sub> alkyl);

R<sup>0a</sup>, R<sup>2a</sup> and R<sup>3a</sup> are each independently -H, -OPg, or halo, wherein Pg is a hydroxy protecting group;

15 R<sup>4</sup> is 1-piperidinyl, 1-pyrrolidinyl, methyl-1-pyrrolidinyl, dimethyl-1-pyrrolidinyl, 4-morpholino, dimethylamino, diethylamino, diisopropylamino, or 1-hexamethyleneimino;

n is 2 or 3;

G<sup>1</sup> is -O-, -S-, or -N(R<sup>5</sup>)-, wherein R<sup>5</sup> is -H or C<sub>1</sub>-C<sub>4</sub> alkyl; and

Y is -O-, -S-, -NH-, -NMe-, or -CH<sub>2</sub>;

20 or a pharmaceutically acceptable salt thereof.

36. A compound according to Claim 35 wherein said compound is 6-(2-hydroxy-benzyl)-5-{hydroxy-[4-(2-piperidin-1-yl-ethoxy)-phenyl]-methyl}-naphthalen-2-ol.